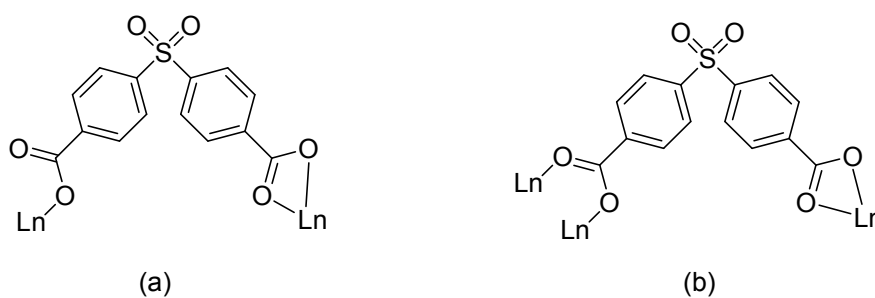


Supporting Information for

A 2D→2D Polyrotaxane Lanthanide–Organic Framework Exhibiting Slow Magnetic Relaxation Behavior

Cai-Ming Liu,* De-Qing Zhang, and Dao-Ben Zhu

*Beijing National Laboratory for Molecular Sciences, Center for Molecular Science,
Institute of Chemistry, Chinese Academy of Sciences Beijing 100190, P.R. China*



Scheme S1. Coordination modes of the L^{2-} ligands.

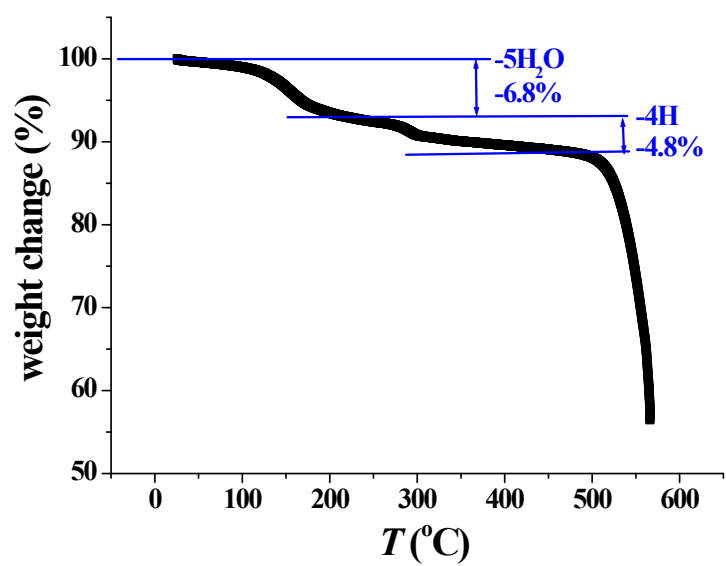


Fig. S1. TGA of **1** measured under a nitrogen atmosphere in the 25-565 °C temperature range and at a scan rate of 10 °Cmin⁻¹.



Figure S2. Coordination geometries of Dy1 and Dy2 in **1**.

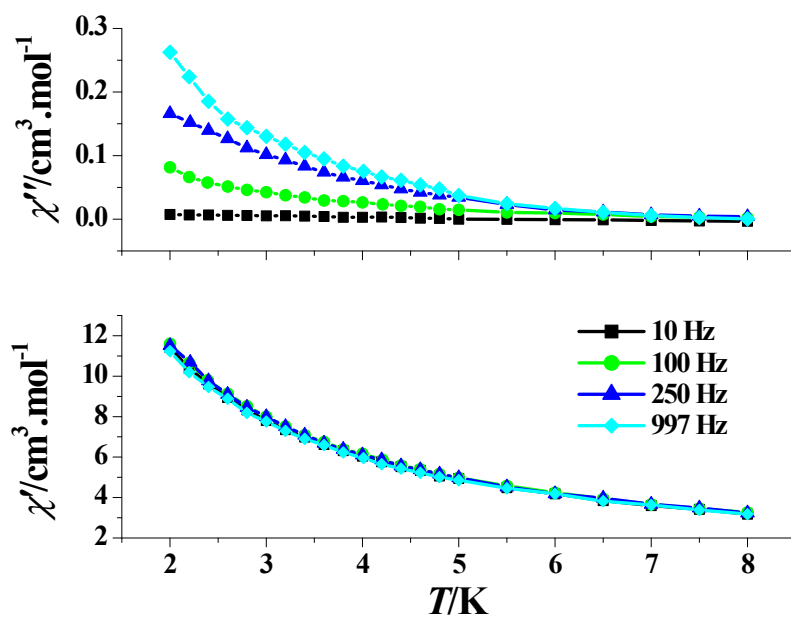


Figure S3. AC susceptibilities measured in a 2.5 Oe ac magnetic field with a zero dc-field for **1**.

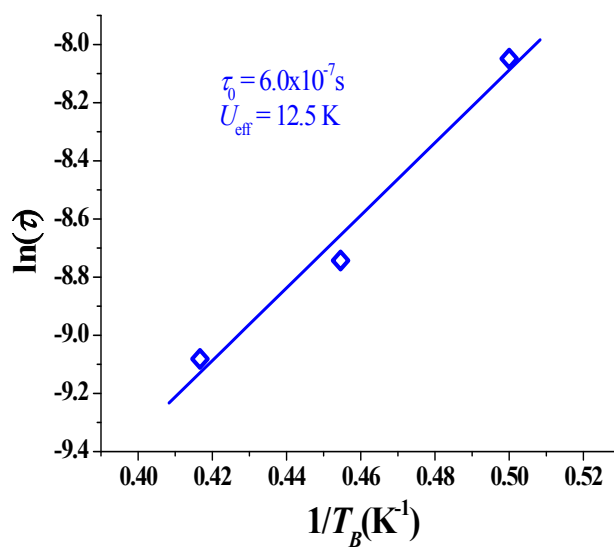


Figure S4. Plot of $\ln(\tau)$ versus $1/T_B$ for **1**, the solid line represents the fitting with the Arrhenius law.

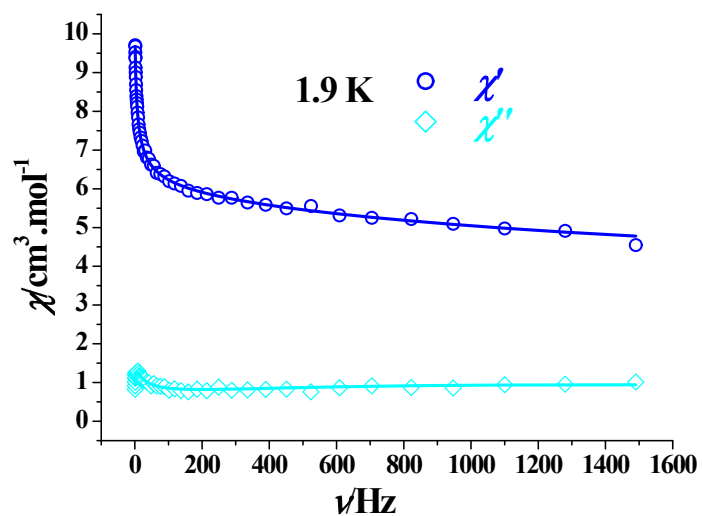


Figure S5. Frequency dependence of the in-phase (χ' , top) and out-of-phase (χ'' , bottom) ac susceptibility of **1** at 1.9 K. the solid lines represent the best fitting with the sum of two modified Debye functions.

Table S1. Linear combination of two modified Debye model fitting parameters at 1.9 K of **1** under 2000 Oe DC field.

$T(\text{K})$	$\chi_2(\text{cm}^3 \cdot \text{mol}^{-1})$	$\chi_1(\text{cm}^3 \cdot \text{mol}^{-1})$	$\chi_0(\text{cm}^3 \cdot \text{mol}^{-1})$	$\tau_1(\text{s})$	α_1	$\tau_2(\text{s})$	α_2
1.9	10.92	8.30	3.14	0.02962	0.435	0.00008	0.305